

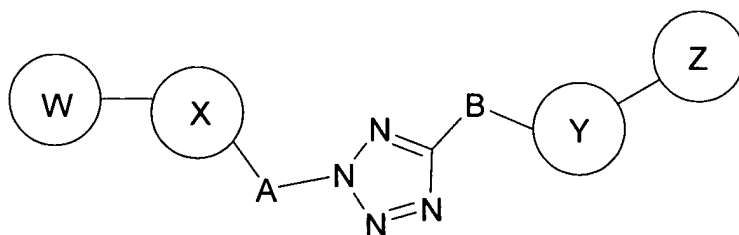
## AMENDMENTS TO THE CLAIMS

Please amend the claims as follows. This listing of claims will replace all prior versions, and listings, of claims in the application.

### Listing of Claims:

Claims 1-37 (Canceled)

38. (New) A compound of the Formula (I):



(I)

wherein:

X is phenyl;

Y is 2-pyridyl;

A is C<sub>0</sub> alkyl,

B is C<sub>0</sub> alkyl,

X is optionally substituted with 1-7 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>1</sup>, -NR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -N(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -NR<sup>1</sup>COR<sup>2</sup>, -NR<sup>1</sup>CO<sub>2</sub>R<sup>2</sup>, -NR<sup>1</sup>SO<sub>2</sub>R<sup>4</sup>, -NR<sup>1</sup>CONR<sup>2</sup>R<sup>3</sup>, -SR<sup>4</sup>, -SOR<sup>4</sup>, -SO<sub>2</sub>R<sup>4</sup>, -SO<sub>2</sub>NR<sup>1</sup>R<sup>2</sup>, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)R<sup>2</sup>, or -C(=NOR<sup>1</sup>)R<sup>2</sup> substituents, wherein optionally two substituents are combined to form a cycloalkyl or heterocycloalkyl ring fused to X; wherein the -C<sub>1-6</sub>alkyl substituent, cycloalkyl ring, or heterocycloalkyl ring each optionally is further substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), or -N(C<sub>0-6</sub>alkyl)(aryl) groups; A is -C<sub>0</sub>alkyl, B is -C<sub>0</sub>alkyl,

Y is optionally substituted with 1-7 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>5</sup>, -NR<sup>5</sup>R<sup>6</sup>, -C(=NR<sup>5</sup>)NR<sup>6</sup>R<sup>7</sup>, -N(=NR<sup>5</sup>)NR<sup>6</sup>R<sup>7</sup>, -NR<sup>5</sup>COR<sup>6</sup>, -NR<sup>5</sup>CO<sub>2</sub>R<sup>6</sup>, -NR<sup>5</sup>SO<sub>2</sub>R<sup>8</sup>, -NR<sup>5</sup>CONR<sup>6</sup>R<sup>7</sup>, -SR<sup>8</sup>, -SOR<sup>8</sup>, -SO<sub>2</sub>R<sup>8</sup>, -SO<sub>2</sub>NR<sup>5</sup>R<sup>6</sup>, -COR<sup>5</sup>, -CO<sub>2</sub>R<sup>5</sup>, -CONR<sup>5</sup>R<sup>6</sup>, -C(=NR<sup>5</sup>)R<sup>6</sup>, or

-C(=NOR<sup>5</sup>)R<sup>6</sup> substituents, wherein optionally two substituents are combined to form a cycloalkyl or heterocycloalkyl ring fused to Y; wherein the -C<sub>1-6</sub>alkyl substituent, cycloalkyl ring, or heterocycloalkyl ring each optionally is further substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), or -N(C<sub>0-6</sub>alkyl)(aryl) groups;

W is -C<sub>3-7</sub>cycloalkyl, -heteroC<sub>3-7</sub>cycloalkyl, -C<sub>0-6</sub>alkylaryl, or -C<sub>0-6</sub>alkylheteroaryl optionally substituted with 1-7 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>1</sup>, -NR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -N(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -NR<sup>1</sup>COR<sup>2</sup>, -NR<sup>1</sup>CO<sub>2</sub>R<sup>2</sup>, -NR<sup>1</sup>SO<sub>2</sub>R<sup>4</sup>, -NR<sup>1</sup>CONR<sup>2</sup>R<sup>3</sup>, -SR<sup>4</sup>, -SOR<sup>4</sup>, -SO<sub>2</sub>R<sup>4</sup>, -SO<sub>2</sub>NR<sup>1</sup>R<sup>2</sup>, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)R<sup>2</sup>, or -C(=NOR<sup>1</sup>)R<sup>2</sup> substituents;

Z is -C<sub>3-7</sub>cycloalkyl, -heteroC<sub>3-7</sub>cycloalkyl, -C<sub>0-6</sub>alkylaryl, or -C<sub>0-6</sub>alkylheteroaryl optionally substituted with 1-7 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>1</sup>, -NR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -N(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -NR<sup>1</sup>COR<sup>2</sup>, -NR<sup>1</sup>CO<sub>2</sub>R<sup>2</sup>, -NR<sup>1</sup>SO<sub>2</sub>R<sup>4</sup>, -NR<sup>1</sup>CONR<sup>2</sup>R<sup>3</sup>, -SR<sup>4</sup>, -SOR<sup>4</sup>, -SO<sub>2</sub>R<sup>4</sup>, -SO<sub>2</sub>NR<sup>1</sup>R<sup>2</sup>, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)R<sup>2</sup>, or -C(=NOR<sup>1</sup>)R<sup>2</sup> substituents;

one of W and Z is optionally absent;

R<sup>1</sup>, R<sup>2</sup>, and R<sup>3</sup> each independently is -C<sub>0-6</sub>alkyl, -C<sub>3-7</sub>cycloalkyl, heteroaryl, or aryl; any of which is optionally substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), -N(C<sub>0-6</sub>alkyl)(aryl) substituents;

R<sup>4</sup> is -C<sub>1-6</sub>alkyl, -C<sub>3-7</sub>cycloalkyl, heteroaryl, or aryl; optionally substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), -N(C<sub>0-6</sub>alkyl)(aryl) substituents;

R<sup>5</sup>, R<sup>6</sup>, and R<sup>7</sup> each independently is -C<sub>0-6</sub>alkyl, -C<sub>3-7</sub>cycloalkyl, heteroaryl, or aryl; any of which is optionally substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), -N(C<sub>0-6</sub>alkyl)(aryl) substituents;

R<sup>8</sup> is -C<sub>1-6</sub>alkyl, -C<sub>3-7</sub>cycloalkyl, heteroaryl, or aryl; optionally substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), -N(C<sub>0-6</sub>alkyl)(aryl) substituents;

R<sup>9</sup> and R<sup>10</sup> each independently is -C<sub>0-6</sub>alkyl, -C<sub>3-7</sub>cycloalkyl, heteroaryl, or aryl; any of which is optionally substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), -N(C<sub>0-6</sub>alkyl)(aryl) substituents; and  
any N may be an N-oxide;  
or a pharmaceutically acceptable salt thereof.

39. (New) The compound according to Claim 38 wherein:

X is phenyl, which is optionally substituted with 1-5 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>1</sup>, -NR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -N(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -NR<sup>1</sup>COR<sup>2</sup>, -NR<sup>1</sup>CO<sub>2</sub>R<sup>2</sup>, -NR<sup>1</sup>SO<sub>2</sub>R<sup>4</sup>, -NR<sup>1</sup>CONR<sup>2</sup>R<sup>3</sup>, -SR<sup>4</sup>, -SOR<sup>4</sup>, -SO<sub>2</sub>R<sup>4</sup>, -SO<sub>2</sub>NR<sup>1</sup>R<sup>2</sup>, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)R<sup>2</sup>, or -C(=NOR<sup>1</sup>)R<sup>2</sup> substituents, wherein optionally two substituents are combined to form a cycloalkyl or heterocycloalkyl ring fused to X; wherein the -C<sub>1-6</sub>alkyl substituent, cycloalkyl ring, or heterocycloalkyl ring each optionally is further substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), or -N(C<sub>0-6</sub>alkyl)(aryl) groups.

40. (New) The compound according to Claim 38 wherein:

Y is 2-pyridyl, which is optionally substituted with 1-4 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>5</sup>, -NR<sup>5</sup>R<sup>6</sup>, -C(=NR<sup>5</sup>)NR<sup>6</sup>R<sup>7</sup>, -N(=NR<sup>5</sup>)NR<sup>6</sup>R<sup>7</sup>, -NR<sup>5</sup>COR<sup>6</sup>, -NR<sup>5</sup>CO<sub>2</sub>R<sup>6</sup>, -NR<sup>5</sup>SO<sub>2</sub>R<sup>8</sup>, -NR<sup>5</sup>CONR<sup>6</sup>R<sup>7</sup>, -SR<sup>8</sup>, -SOR<sup>8</sup>, -SO<sub>2</sub>R<sup>8</sup>, -SO<sub>2</sub>NR<sup>5</sup>R<sup>6</sup>, -COR<sup>5</sup>, -CO<sub>2</sub>R<sup>5</sup>, -CONR<sup>5</sup>R<sup>6</sup>, -C(=NR<sup>5</sup>)R<sup>6</sup>, or -C(=NOR<sup>5</sup>)R<sup>6</sup> substituents, wherein optionally two substituents are combined to form a cycloalkyl or heterocycloalkyl ring fused to Y; wherein the -C<sub>1-6</sub>alkyl substituent, cycloalkyl ring, or heterocycloalkyl ring each optionally is further substituted with 1-5 independent halogen, -CN, -C<sub>1-6</sub>alkyl, -O(C<sub>0-6</sub>alkyl), -O(C<sub>3-7</sub>cycloalkyl), -O(aryl), -N(C<sub>0-6</sub>alkyl)(C<sub>0-6</sub>alkyl), -N(C<sub>0-6</sub>alkyl)(C<sub>3-7</sub>cycloalkyl), or -N(C<sub>0-6</sub>alkyl)(aryl) groups.

41. (New) The compound according to Claim 38 wherein:

W is -C<sub>0-6</sub>alkylaryl, or -C<sub>0-6</sub>alkylheteroaryl optionally substituted with 1-7 independent halogen, -CN, NO<sub>2</sub>, -C<sub>1-6</sub>alkyl, -C<sub>1-6</sub>alkenyl, -C<sub>1-6</sub>alkynyl, -OR<sup>1</sup>, -NR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -N(=NR<sup>1</sup>)NR<sup>2</sup>R<sup>3</sup>, -NR<sup>1</sup>COR<sup>2</sup>, -NR<sup>1</sup>CO<sub>2</sub>R<sup>2</sup>, -NR<sup>1</sup>SO<sub>2</sub>R<sup>4</sup>, -NR<sup>1</sup>CONR<sup>2</sup>R<sup>3</sup>, -SR<sup>4</sup>, -SOR<sup>4</sup>, -SO<sub>2</sub>R<sup>4</sup>, -SO<sub>2</sub>NR<sup>1</sup>R<sup>2</sup>, -COR<sup>1</sup>, -CO<sub>2</sub>R<sup>1</sup>, -CONR<sup>1</sup>R<sup>2</sup>, -C(=NR<sup>1</sup>)R<sup>2</sup>, or -C(=NOR<sup>1</sup>)R<sup>2</sup> substituents.

42. (New) A compound which is selected from the group consisting of:

1-methyl-3-[3-(5-pyridin-2-yl-2*H*-tetrazol-2-yl)phenyl]imidazolidin-2-one;

2-[2-(4-pyridin-2-ylphenyl)-2*H*-tetrazol-5-yl]pyridine;

2-[2-(4-pyridin-4-ylphenyl)-2*H*-tetrazol-5-yl]pyridine;

2-{2-[3-(1*H*-imidazol-1-yl)phenyl]-2*H*-tetrazol-5-yl}pyridine;

2-[2-(2-pyrazin-3-ylphenyl)-2*H*-tetrazol-5-yl]pyridine;

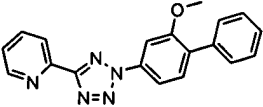
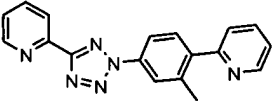
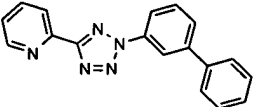
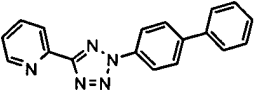
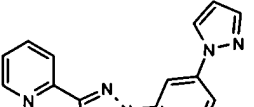
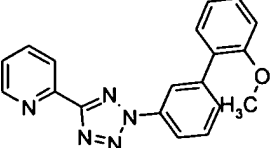
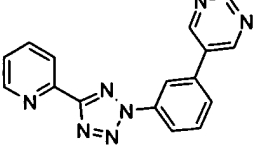
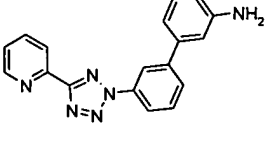
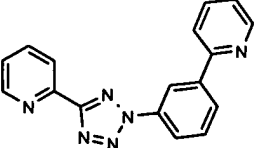
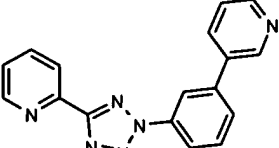
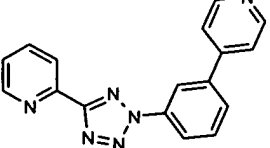
2-[2-(4-morpholin-3-ylphenyl)-2*H*-tetrazol-5-yl]pyridine;

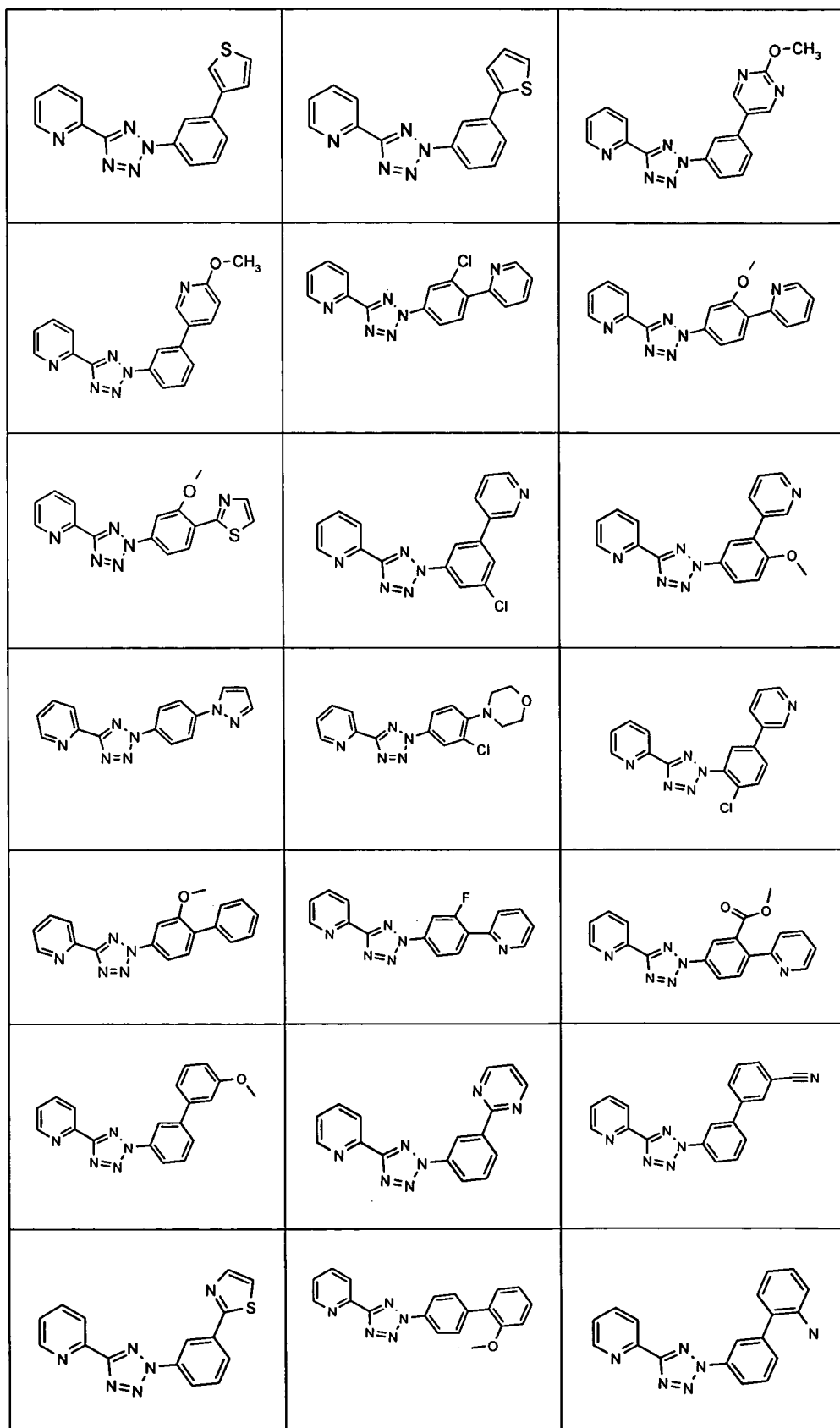
2-{2-[3-(2*H*-tetrazol-5-yl)phenyl]-2*H*-tetrazol-5-yl}pyridine; and

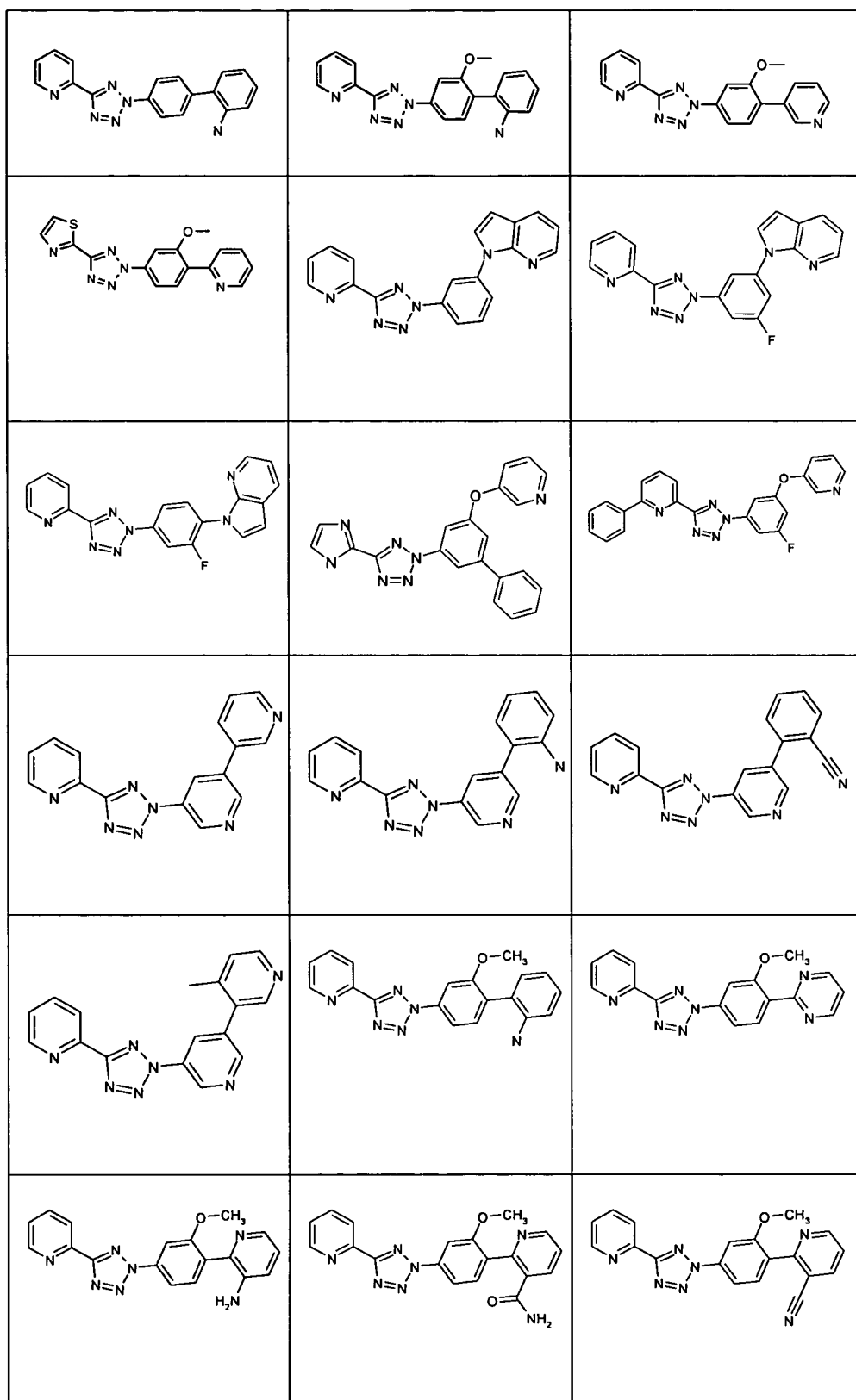
2-pyridin-2-yl-5-(5-pyridin-2-yl-2*H*-tetrazol-2-yl)benzonitrile;

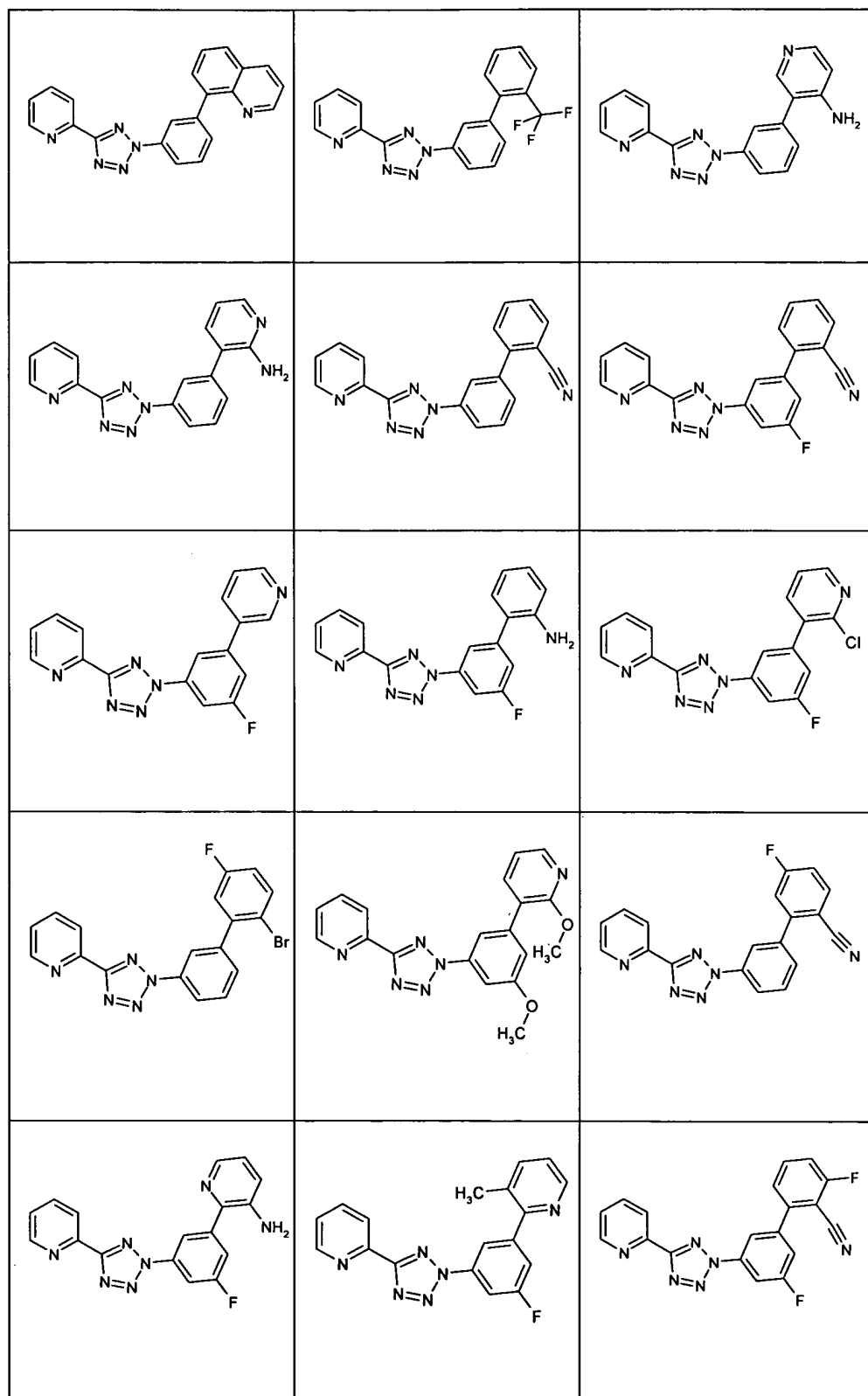
or a pharmaceutically acceptable salt thereof.

43. (New) A compound which selected from the group consisting of:

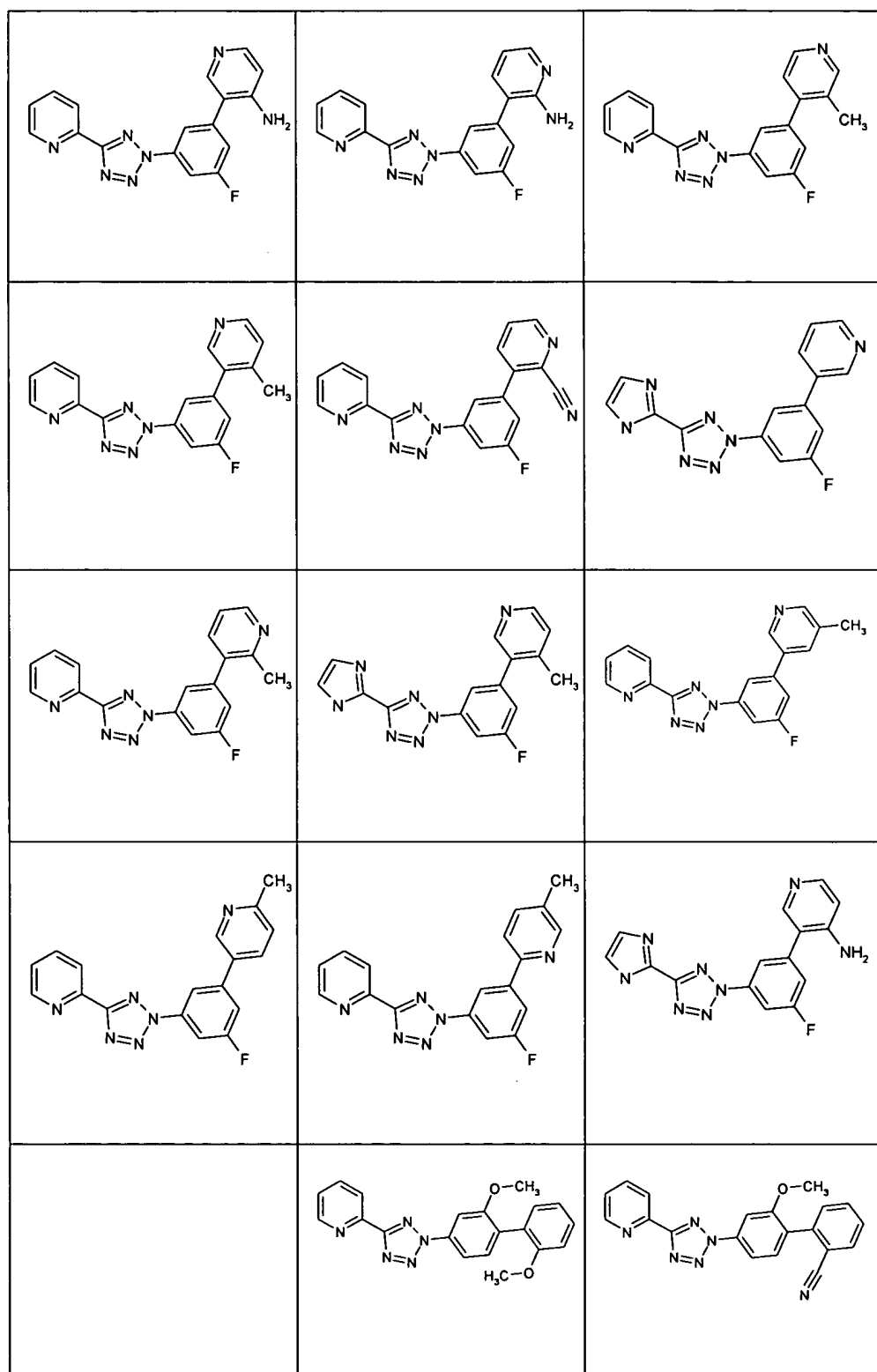
		
		
		
		

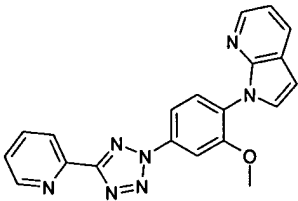
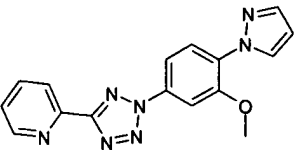
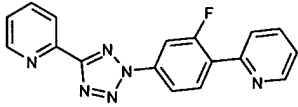
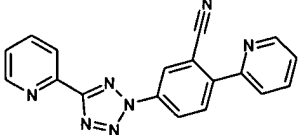
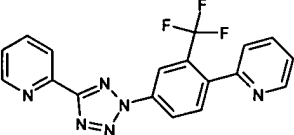
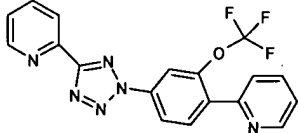
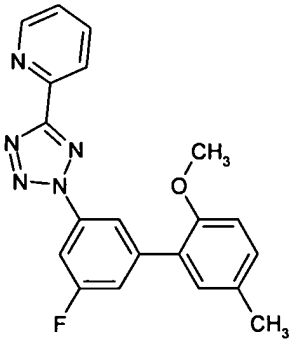
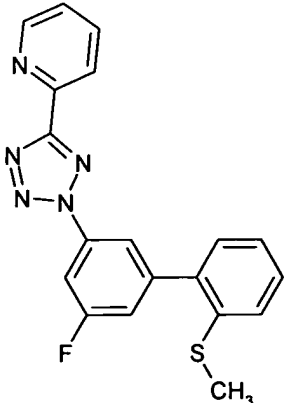


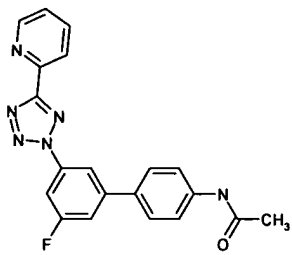
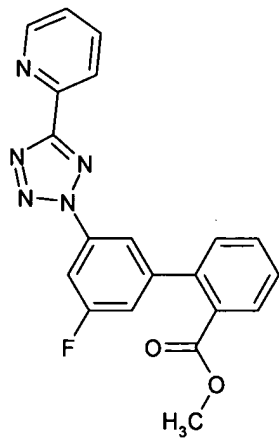
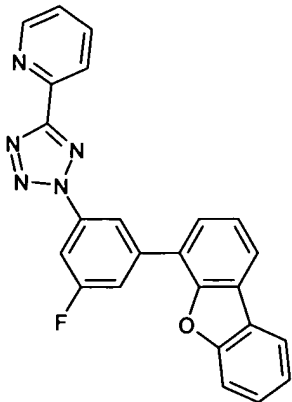


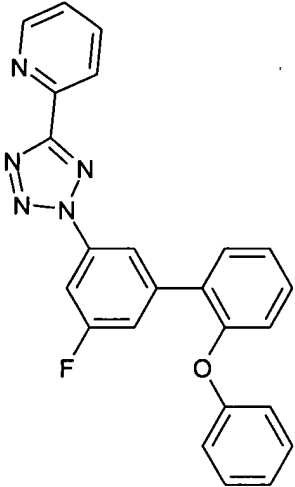
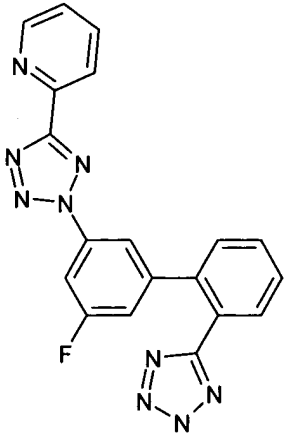
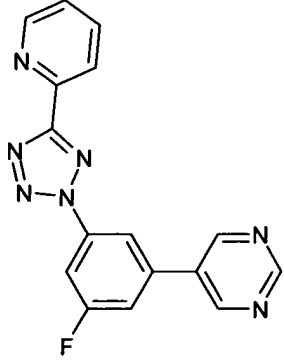



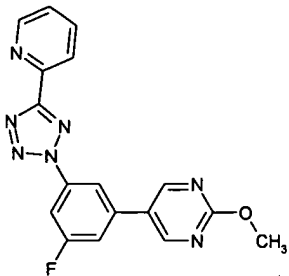
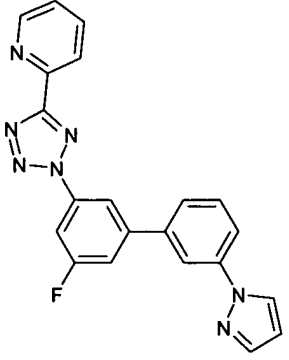
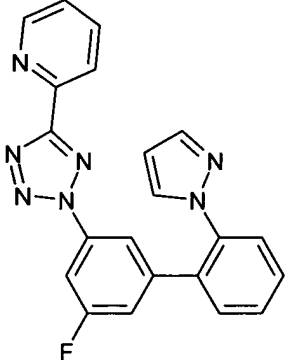
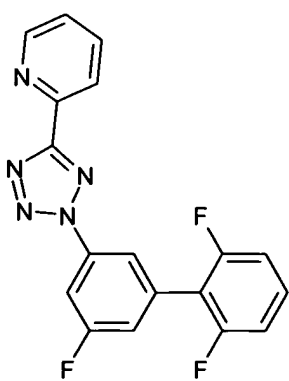


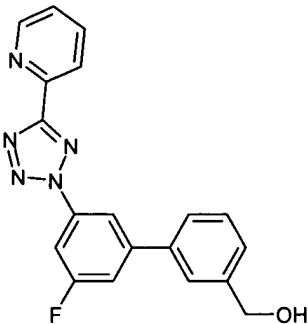
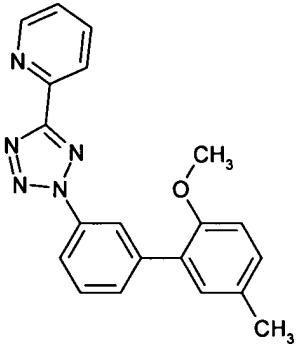
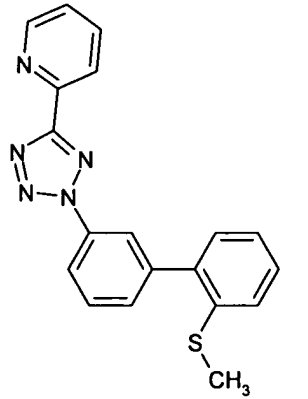


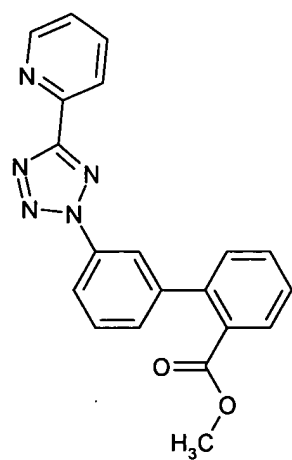
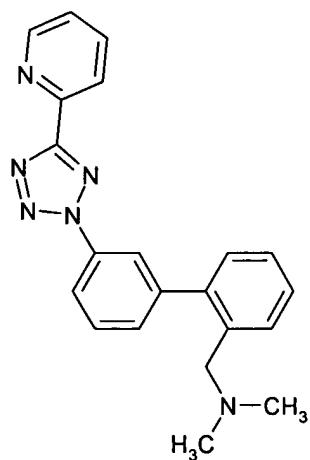
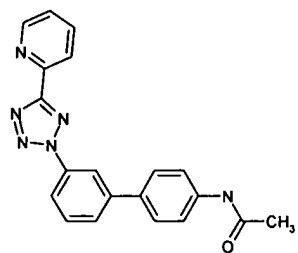
		
		
		
		
		

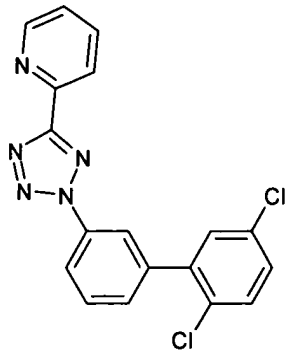
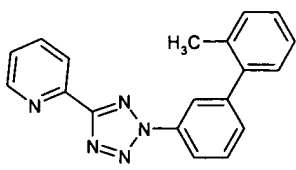
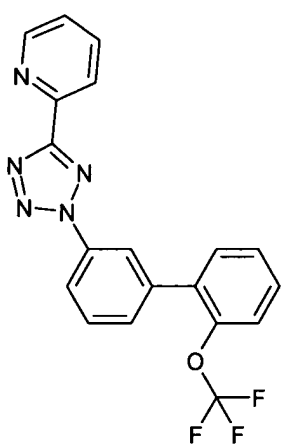
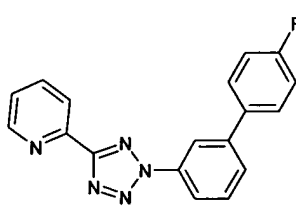
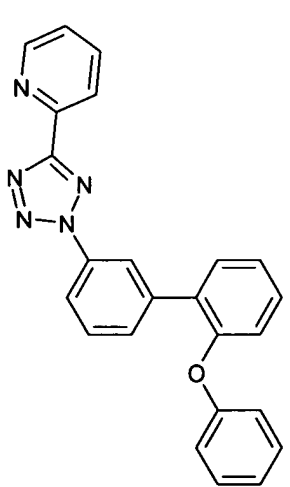
		
		
		

	 <p>Chemical structure of 1-(4-(4-(benzyloxy)phenyl)-5-(pyridin-2-yl)-1H-1,2,4-triazol-3-yl)benzene. The structure features a central benzene ring substituted at the 1-position with a 1,2,4-triazole ring, at the 3-position with a pyridin-2-yl group, and at the 4-position with a 4-(benzyloxy)phenyl group.</p>	
	 <p>Chemical structure of 1-(4-(4-(1,2,4-triazol-5-yl)phenyl)-5-(pyridin-2-yl)-1H-1,2,4-triazol-3-yl)benzene. The structure features a central benzene ring substituted at the 1-position with a 1,2,4-triazole ring, at the 3-position with a pyridin-2-yl group, and at the 4-position with a 4-(1,2,4-triazol-5-yl)phenyl group.</p>	
	 <p>Chemical structure of 1-(4-(4-(pyridin-2-yl)-5-(1,2,4-triazol-5-yl)phenyl)-5-(pyridin-2-yl)-1H-1,2,4-triazol-3-yl)benzene. The structure features a central benzene ring substituted at the 1-position with a 1,2,4-triazole ring, at the 3-position with a pyridin-2-yl group, and at the 4-position with a 4-(pyridin-2-yl)-5-(1,2,4-triazol-5-yl)phenyl group.</p>	

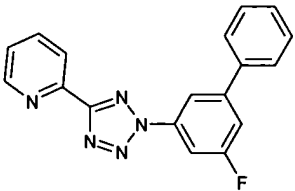
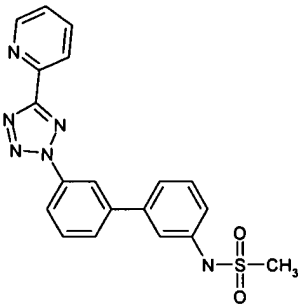
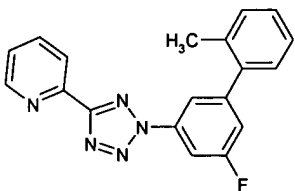
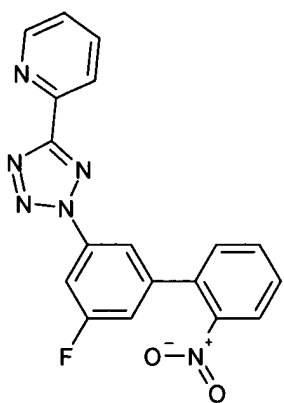
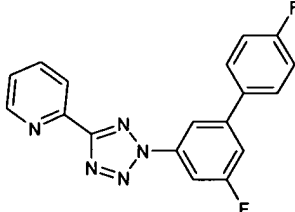
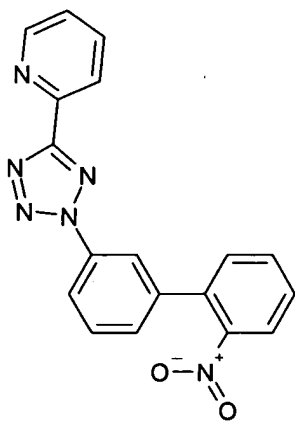
		
		
		
		

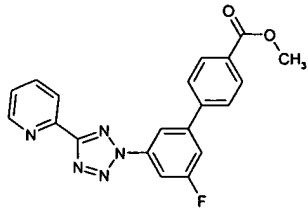
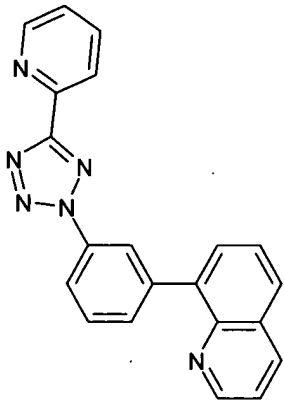
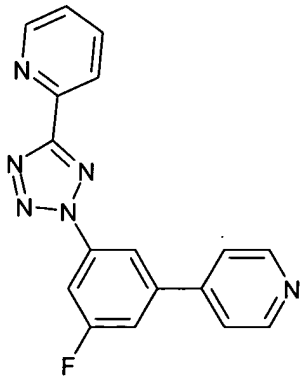
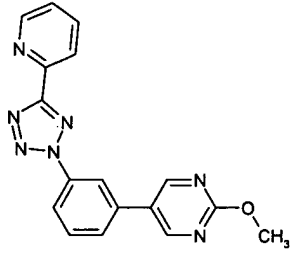
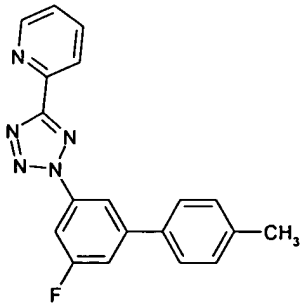
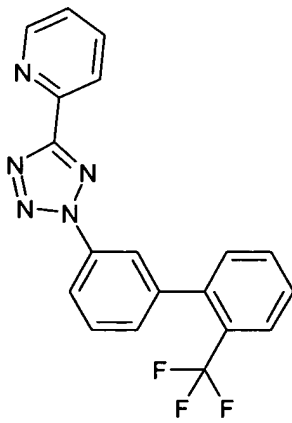
	 <p>Chemical structure of 4-(4-(benzo[d][1,2,4]triazol-1-yl)phenyl)-2-fluorophenylmethanol. It features a central biphenyl core. The left ring is substituted with a fluorine atom at the 2-position and a 1,2,4-benzotriazol-1-yl group at the 4-position. The right ring is substituted with a hydroxymethyl group (-CH<sub>2</sub>OH) at the 4-position.</p>	
	 <p>Chemical structure of 4-(4-(benzo[d][1,2,4]triazol-1-yl)phenyl)-2-methoxyphenylmethane. It features a central biphenyl core. The left ring is substituted with a 1,2,4-benzotriazol-1-yl group at the 4-position. The right ring is substituted with a methoxy group (-OCH<sub>3</sub>) at the 2-position and a methyl group (-CH<sub>3</sub>) at the 4-position.</p>	
	 <p>Chemical structure of 4-(4-(benzo[d][1,2,4]triazol-1-yl)phenyl)-2-(methylsulfanyl)phenylmethane. It features a central biphenyl core. The left ring is substituted with a 1,2,4-benzotriazol-1-yl group at the 4-position. The right ring is substituted with a methylsulfanyl group (-SCH<sub>3</sub>) at the 2-position.</p>	

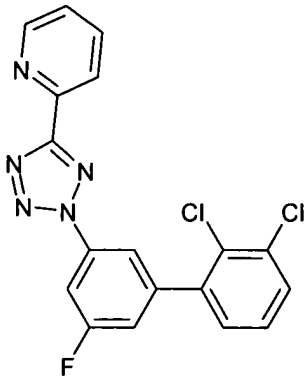
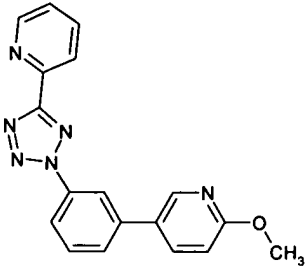
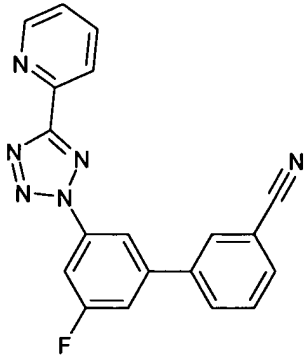
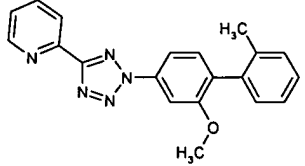
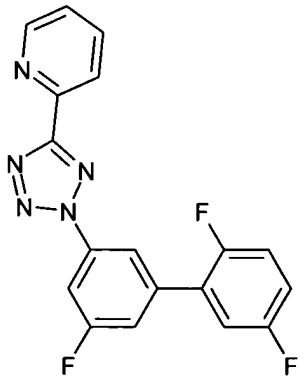
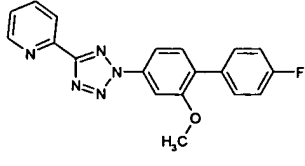


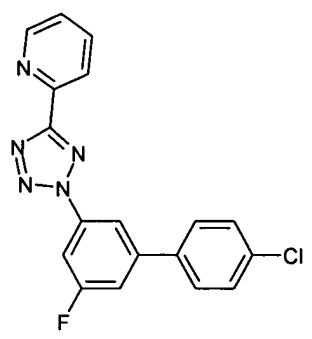
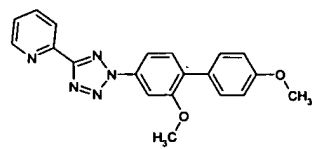
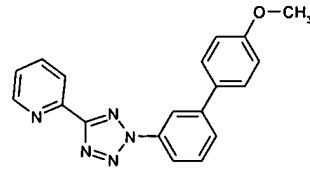
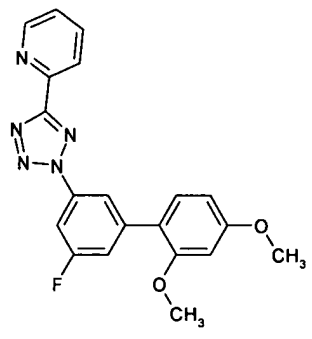
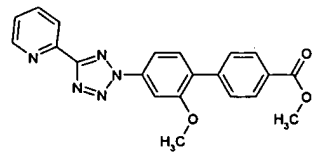
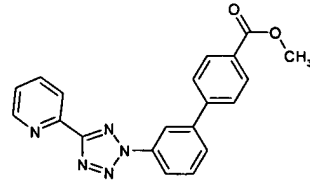
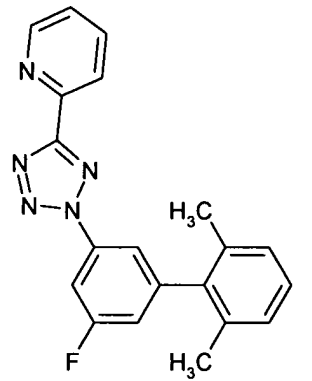
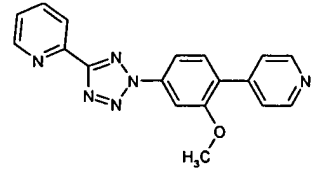
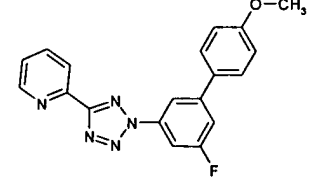
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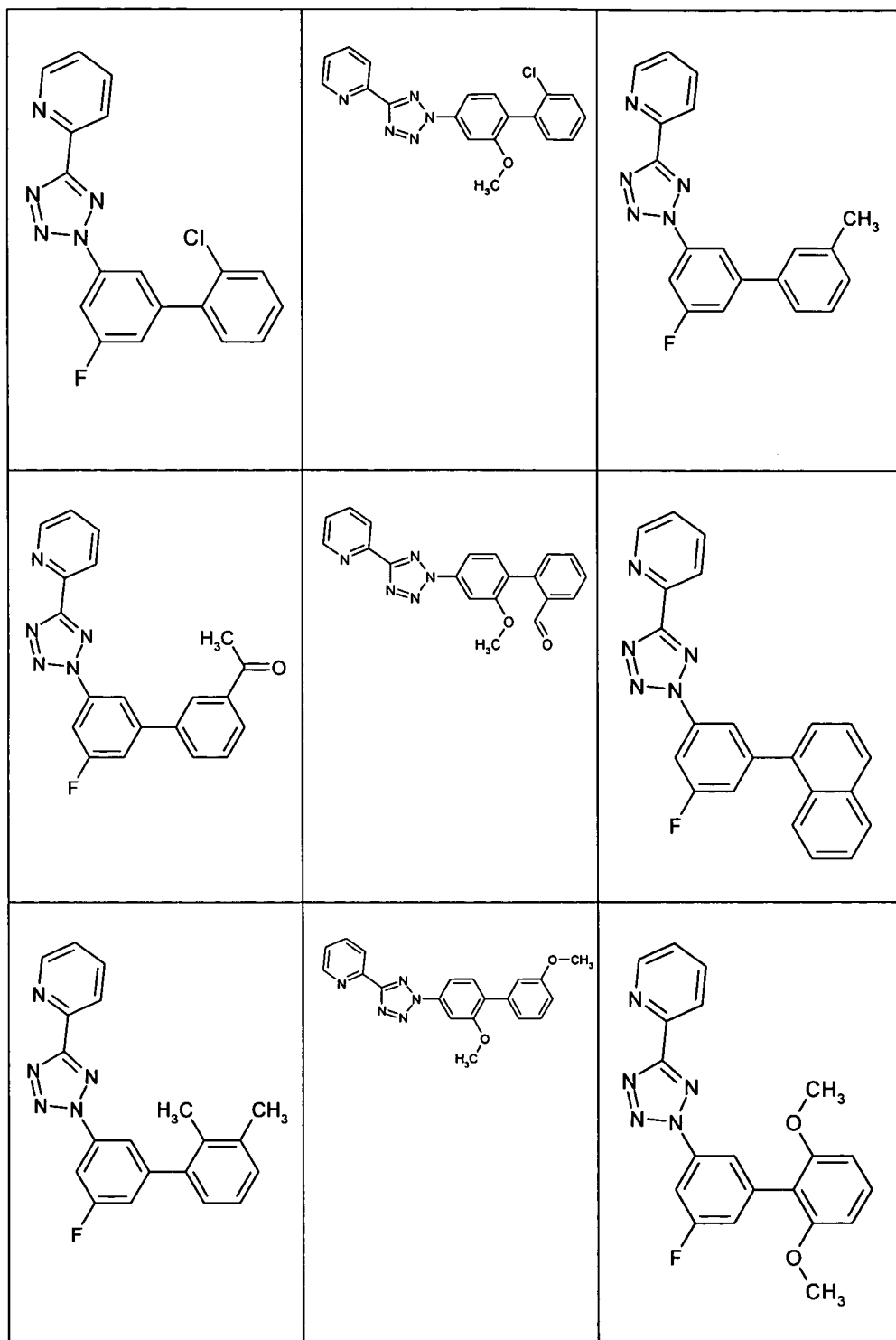


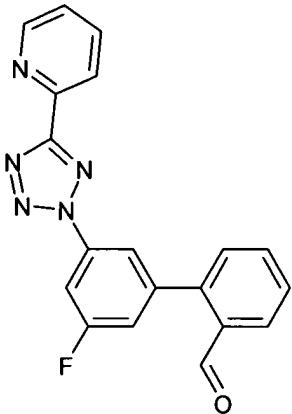
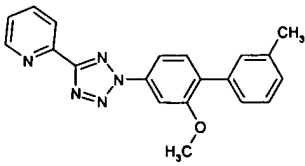
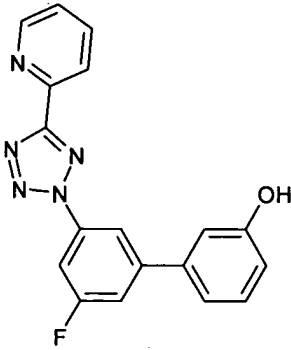
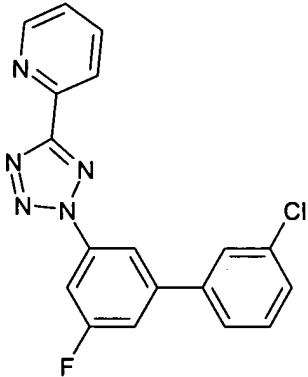
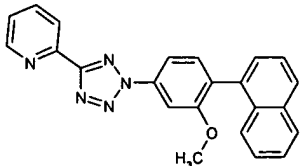
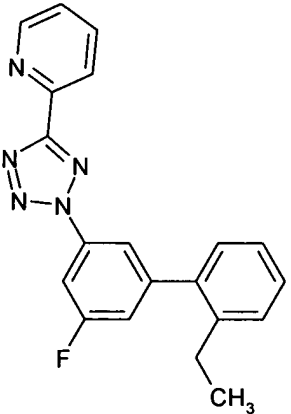
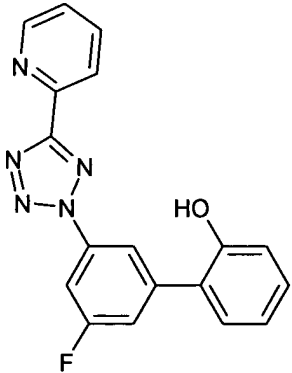
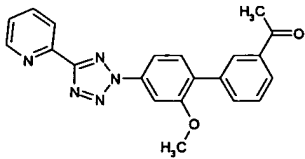
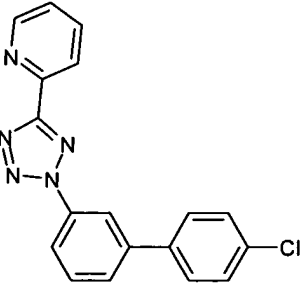
		
		
		

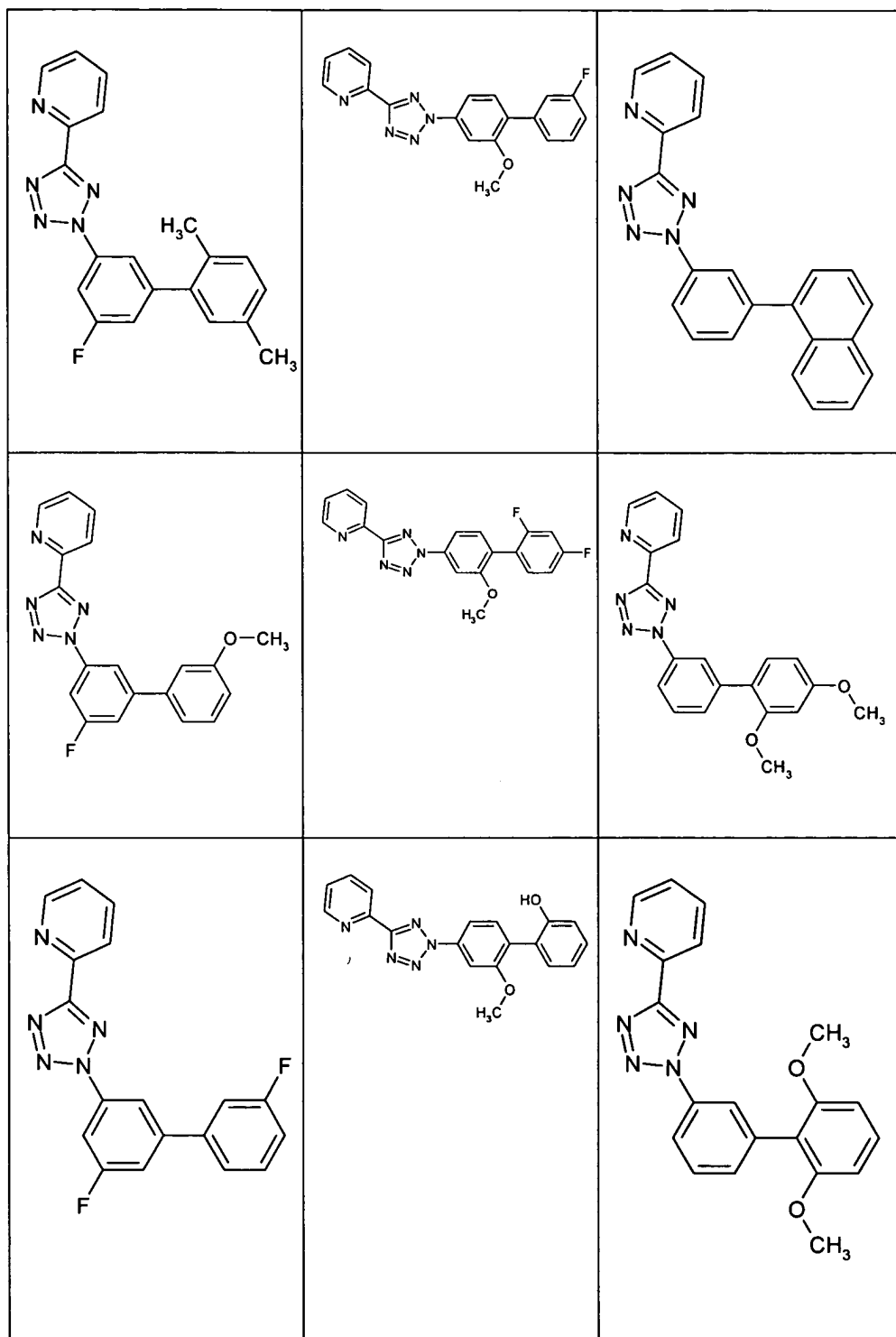
		
		
		

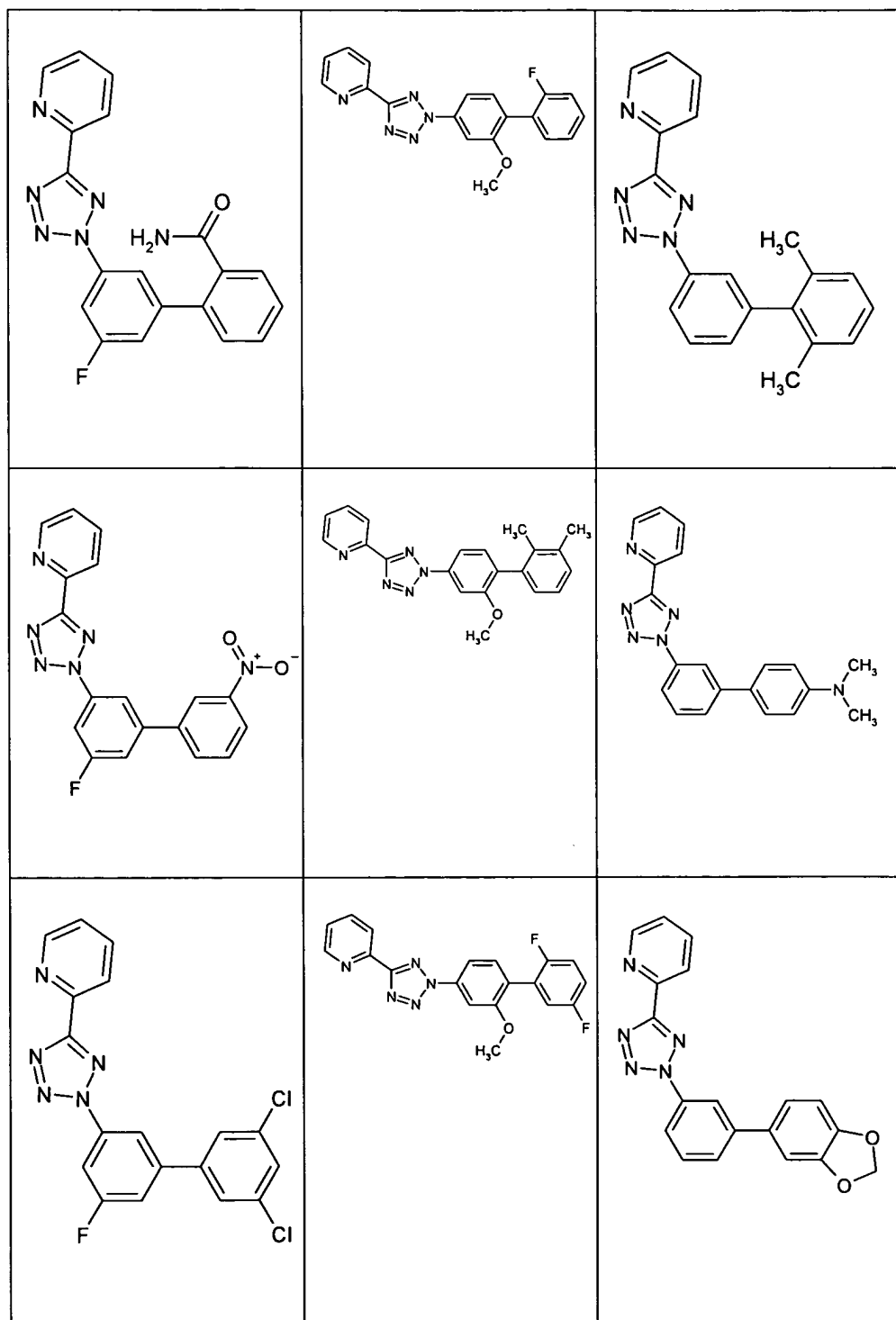
		
		
		

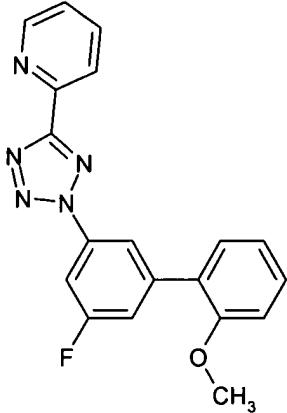
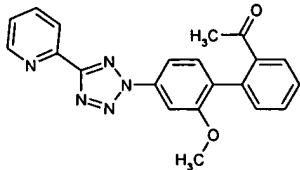
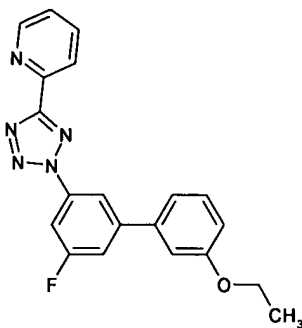
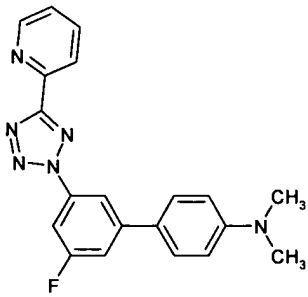
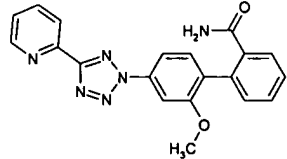
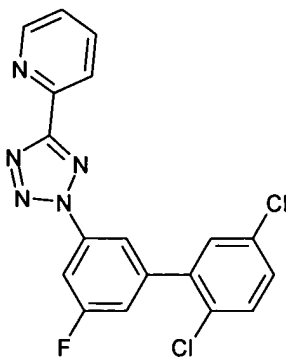
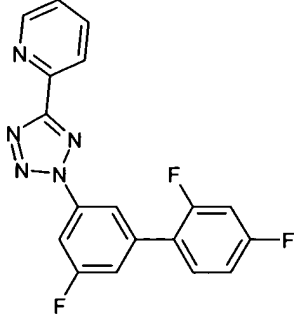
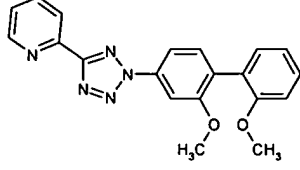
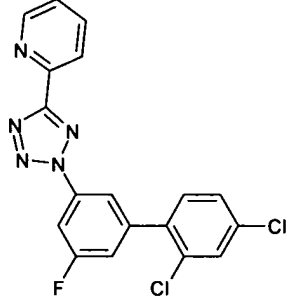


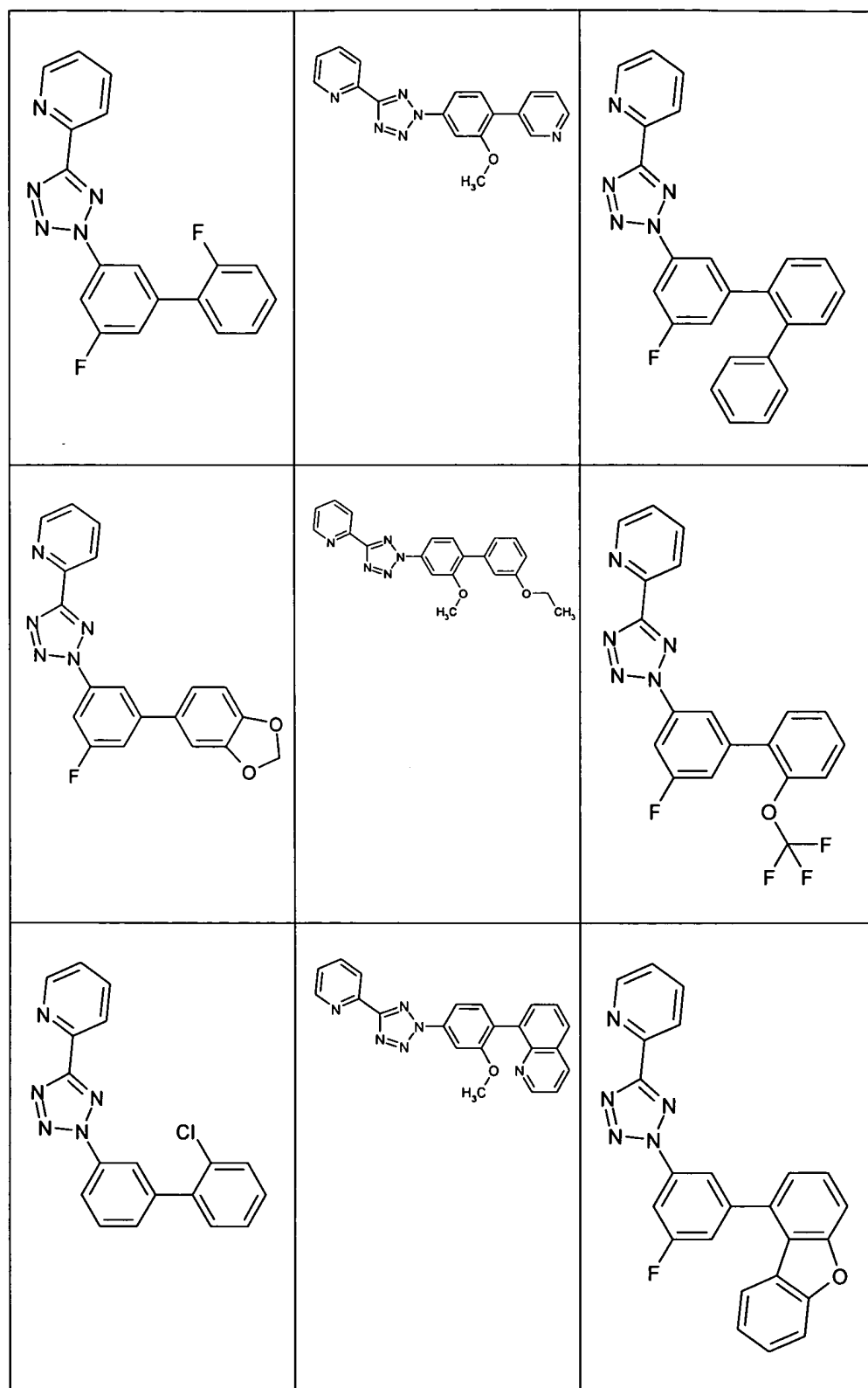
		
		
		

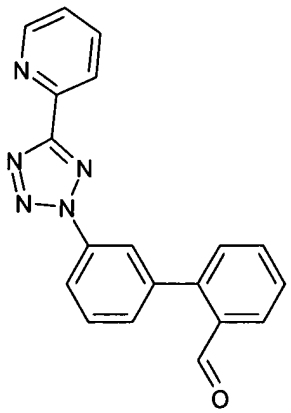
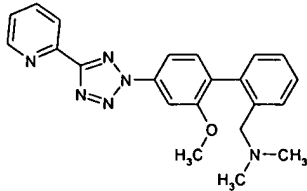
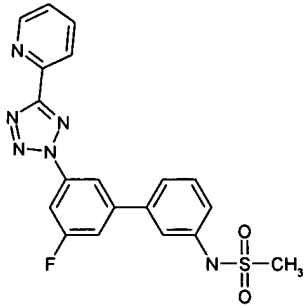
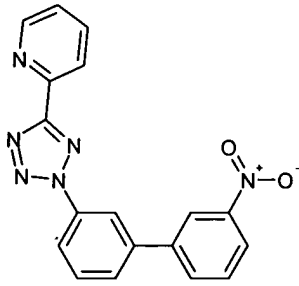
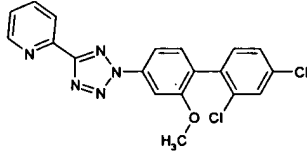
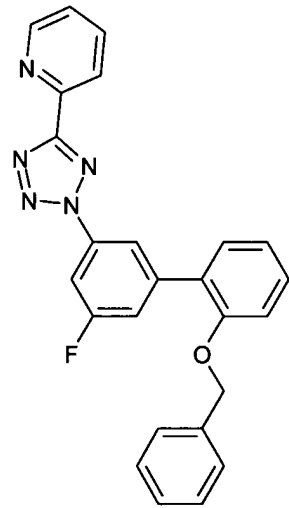
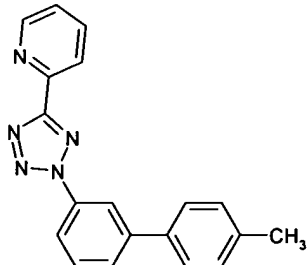
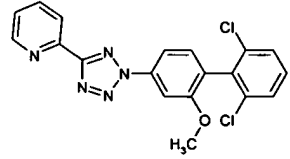
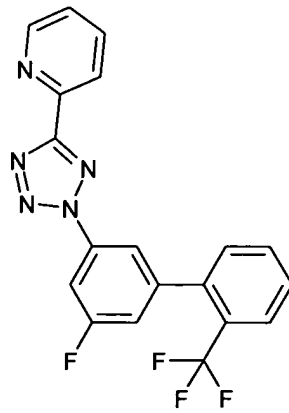


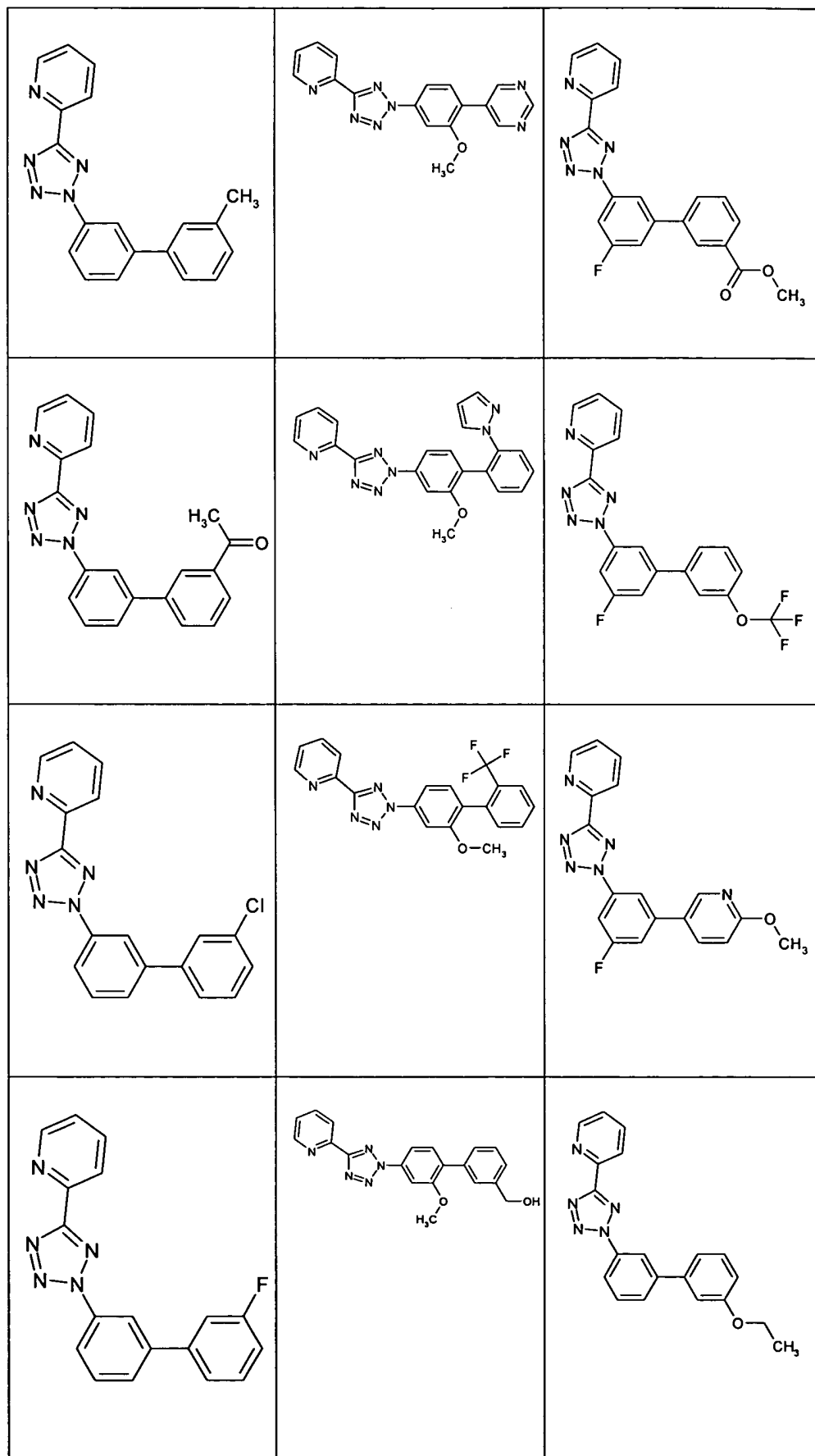


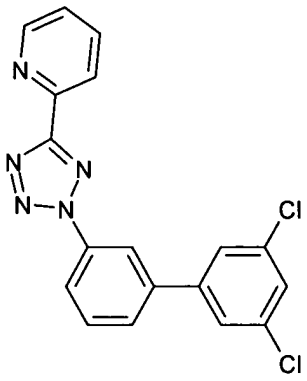
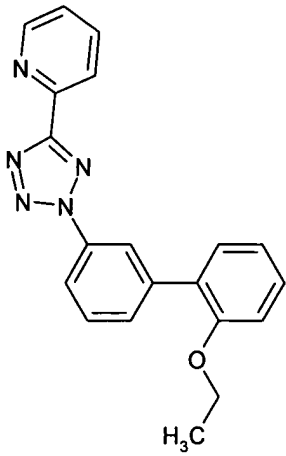
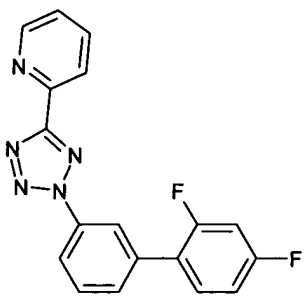
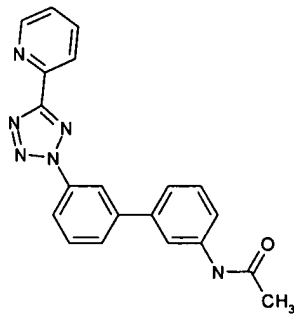
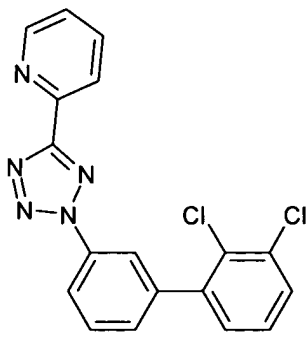
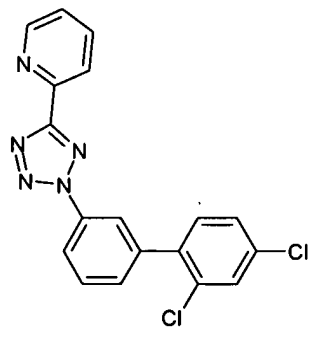


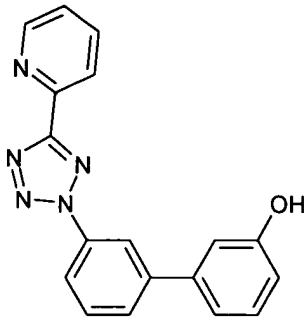
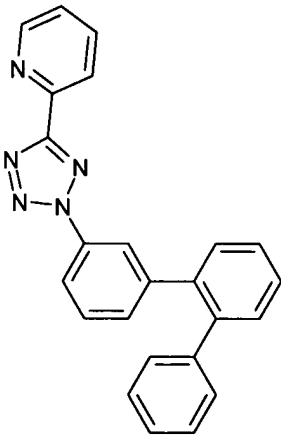
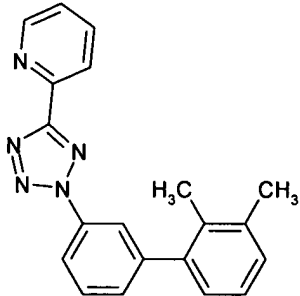
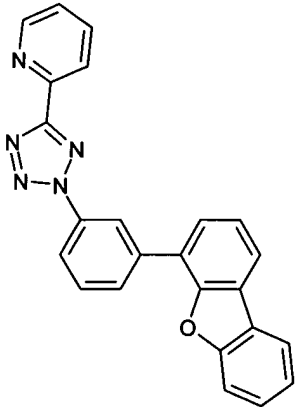
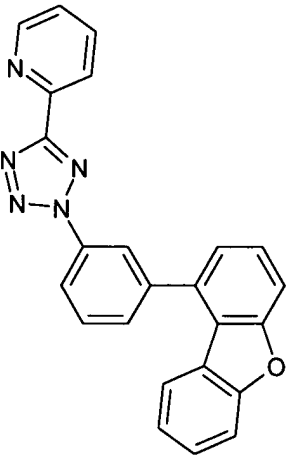
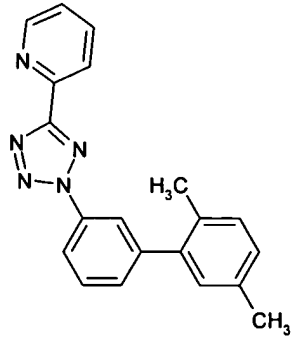
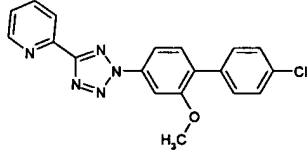
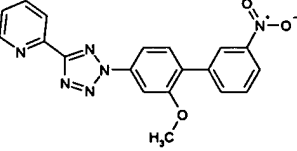
		
		
		

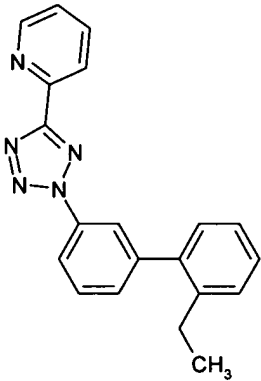
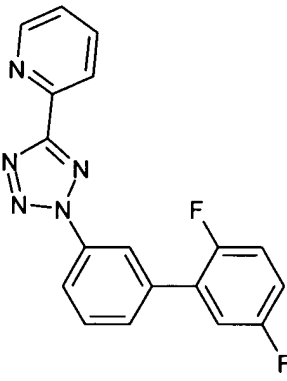
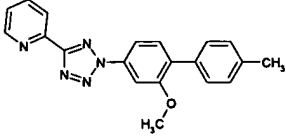
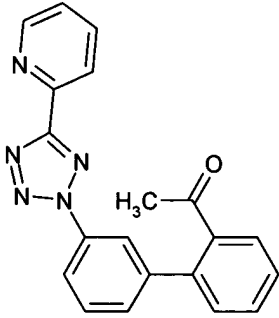
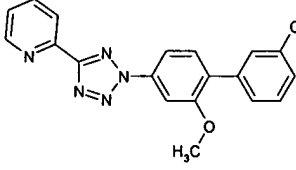
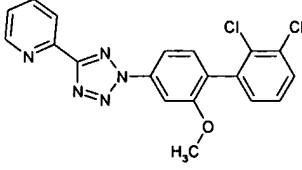
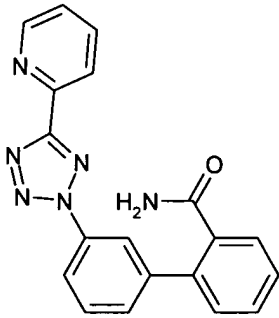
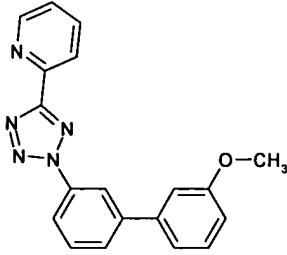
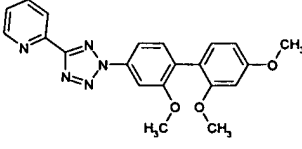
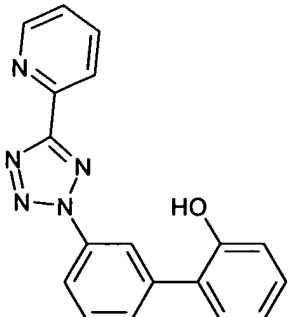
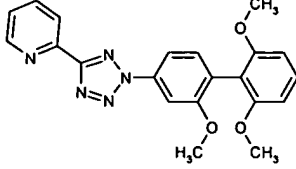
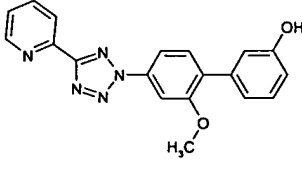


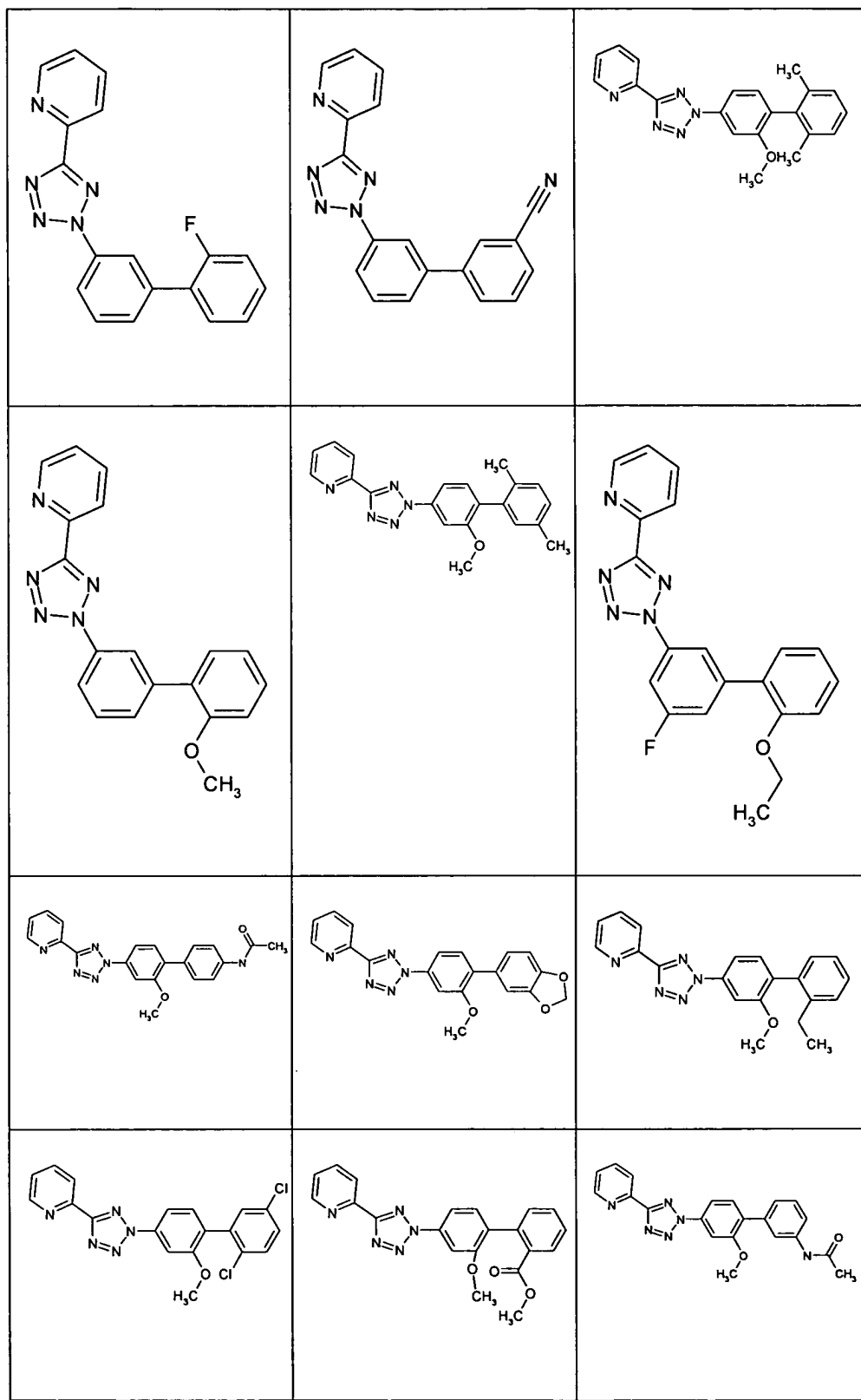
		
		
		



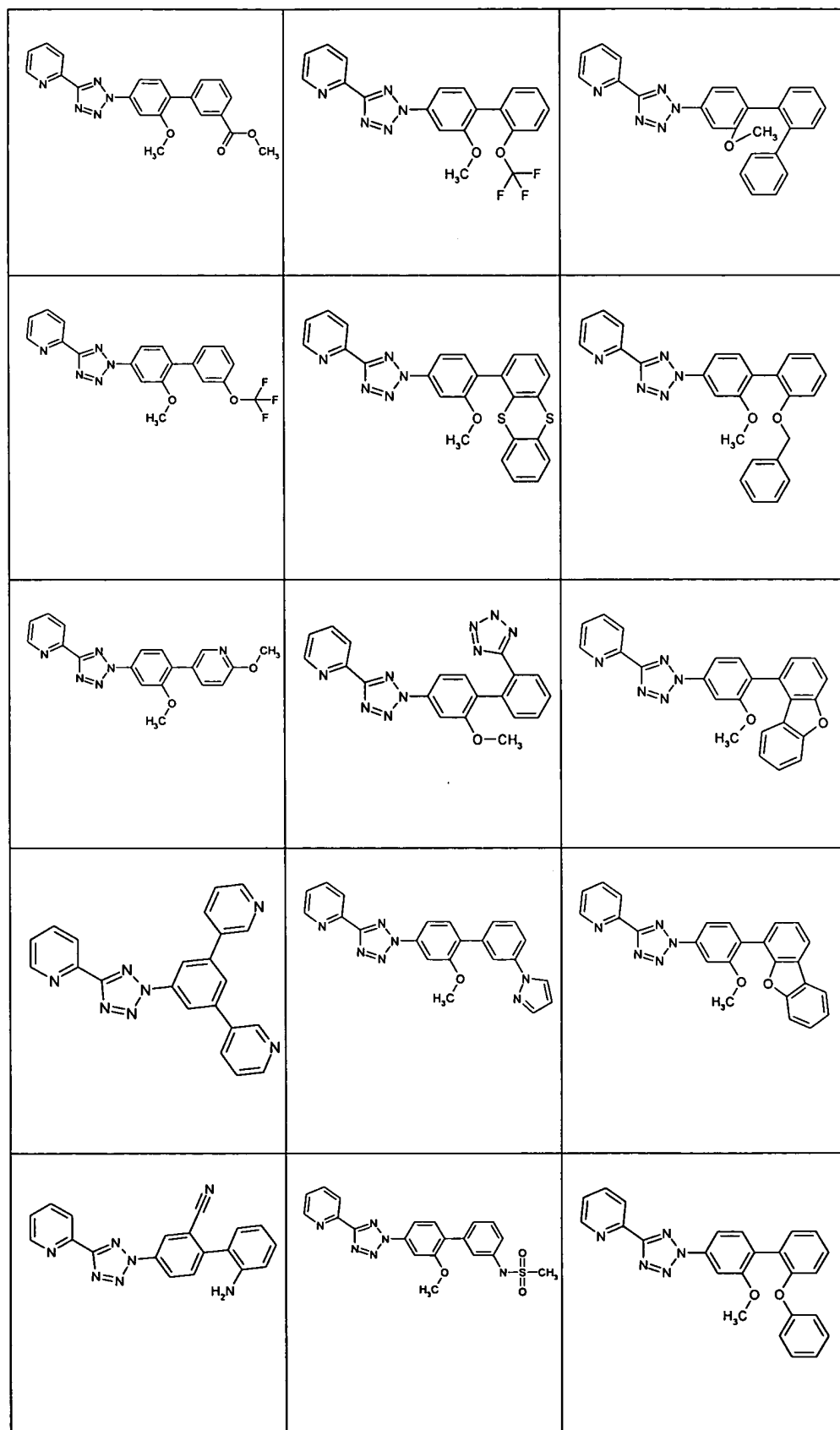
 <chem>Clc1ccc(cc1-c2ccccc2-n3nn[nH]3-c4ccccn4)Cl</chem>		 <chem>CCOC1=CC=C(C=C1-c2ccccc2-n3nn[nH]3-c4ccccn4)</chem>
 <chem>Fc1cc(F)ccc1-c2ccccc2-n3nn[nH]3-c4ccccn4</chem>		 <chem>CC(=O)Nc1ccc(cc1-c2ccccc2-n3nn[nH]3-c4ccccn4)</chem>
 <chem>Clc1cc(Cl)ccc1-c2ccccc2-n3nn[nH]3-c4ccccn4</chem>		 <chem>Clc1cc(Cl)ccc1-c2ccccc2-n3nn[nH]3-c4ccccn4</chem>

 <chem>Oc1ccc(cc1)-c2ccc(cc2)-n3nn[nH]3c4ccncc4</chem>		 <chem>c1ccc(cc1)-c2ccccc2-c3ccc(cc3)-c4ccc(cc4)-n5nn[nH]5c6ccncc6</chem>
 <chem>Cc1cc(C)ccc1-c2ccc(cc2)-n3nn[nH]3c4ccncc4</chem>	 <chem>c1ccc2c(c1)oc3ccccc3c2-c4ccc(cc4)-n5nn[nH]5c6ccncc6</chem>	 <chem>c1ccc2c(c1)oc3ccccc3c2-c4ccc(cc4)-n5nn[nH]5c6ccncc6</chem>
 <chem>Cc1cc(C)ccc1-c2ccc(cc2)-n3nn[nH]3c4ccncc4</chem>	 <chem>COc1cc(ccc1-c2ccc(cc2)-n3nn[nH]3c4ccncc4)-c5ccc(Cl)cc5</chem>	 <chem>[O-][N+](=O)c1ccc(cc1-c2ccc(cc2)-n3nn[nH]3c4ccncc4)-c5cc(OC)cc5</chem>







or a pharmaceutically acceptable salt thereof.

44. (New) A pharmaceutical composition comprising the compound of Claim 38, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

45. (New) A pharmaceutical composition comprising the compound of Claim 42, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

46. (New) A pharmaceutical composition comprising the compound of Claim 43, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.